

EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	115	562/586	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 07:21
L2	1572	((562/586) or (562/856) or (568/486) or (568/488) or (568/489) or (568/490) or (568/615) or (568/622)). CCLS.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	OFF	2006/08/31 08:45
L3	8204727	reduc\$5 or hydrogen\$6	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 07:24
L4	964	hydrofluoroether	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 07:25
L5	1940726	platinum or pt!	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 08:45
L6	232	(I2 or I4) and I3 and I5	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 07:26
L7	81722	catalyst near10 I5	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 07:28
L8	81	I6 and I7	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 08:40
L9	660	Picozzi.in. or Meo.in. or Tonelli.in.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 08:41

EAST Search History

L10	7	I3 and I5 and I7 and I9	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 08:45
L11	1454	((562/586) or (562/856) or (568/486) or (568/488) or (568/489) or (568/490) or (568/615) or (568/622)). CCLS.	US-PGPUB; USPAT; USOCR	OR	OFF	2006/08/31 08:45
L12	35047	platinum.clm.	US-PGPUB; USPAT; USOCR	OR	ON	2006/08/31 08:46
L13	16	I11 and I12	US-PGPUB; USPAT; USOCR	OR	ON	2006/08/31 08:46

10/630,697

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NEWS 10 JUN 02 The first reclassification of IPC codes now complete in
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NEWS 11 JUN 26 TULSA/TULSA2 reloaded and enhanced with new search and
and display fields
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NEWS 13 JUL 11 CHEMSAFE reloaded and enhanced
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NEWS 15 JUL 19 Coverage of Research Disclosure reinstated in DWPI
NEWS 16 AUG 09 INSPEC enhanced with 1898-1968 archive
NEWS 17 AUG 28 ADISCTI Reloaded and Enhanced
NEWS 18 AUG 30 CA(SM)/CAPLUS(SM) Austrian patent law changes

NEWS EXPRESS JUNE 30 CURRENT WINDOWS VERSION IS V8.01b, CURRENT
MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
AND CURRENT DISCOVER FILE IS DATED 26 JUNE 2006.

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FILE 'HOME' ENTERED AT 07:02:32 ON 31 AUG 2006

=> file reg

COST IN U.S. DOLLARS

SINCE FILE
ENTRY

TOTAL
SESSION

10/630,697

FULL ESTIMATED COST

0.21

0.21

FILE 'REGISTRY' ENTERED AT 07:02:59 ON 31 AUG 2006
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=>Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=> screen 963

L1 SCREEN CREATED

=> screen 1992 OR 2016 OR 2021 OR 2026 OR 1838

L2 SCREEN CREATED

=>

Uploading C:\Documents and Settings\rkeys\My
Documents\STNEXP4\TEMPLATE\STANDARD\10630697.str



10/630,697

chain nodes :

1 2 3 4 5 6 7 9 10 11 12 13 15 16 17 18

chain bonds :

1-2 2-3 2-4 4-5 4-6 4-9 6-7 7-18 10-11 11-12 11-13 11-15 15-16 16-17

exact/norm bonds :

2-3 4-6 4-9 6-7 11-13 11-15 15-16

exact bonds :

1-2 2-4 4-5 7-18 10-11 11-12 16-17

G1:F,CF3

Match level :

1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 9:CLASS 10:CLASS

11:CLASS 12:CLASS 13:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS

fragments assigned product role:

containing 10

fragments assigned reactant/reagent role:

containing 1

L3 STRUCTURE UPLOADED

=> que L3 AND L1 NOT L2

L4 QUE L3 AND L1 NOT L2

=> d

L4 HAS NO ANSWERS

L1 SCR 963

L2 SCR 1992 OR 2016 OR 2021 OR 2026 OR 1838

L3 STR

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.

L4 QUE L3 AND L1 NOT L2

=> file reaction

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

0.88

1.09

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=> s 14

SAMPLE SEARCH INITIATED 07:04:11 FILE 'CASREACT'

SCREENING COMPLETE - 8 REACTIONS TO VERIFY FROM

2 DOCUMENTS

100.0% DONE

8 VERIFIED

0 HIT RXNS

0 DOCS

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SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED VERIFICATIONS: 8 TO 329
PROJECTED ANSWERS: 0 TO 0

SAMPLE SEARCH INITIATED 07:04:13 FILE 'CHEMINFORMRX'
SCREENING COMPLETE - 3 REACTIONS TO VERIFY FROM 1 DOCUMENTS

100.0% DONE 3 VERIFIED 0 HIT RXNS 0 DOCS
SEARCH TIME: 00.00.07

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED VERIFICATIONS: 3 TO 163
PROJECTED ANSWERS: 0 TO 0

FULL SEARCH INITIATED 07:04:23 FILE 'DJSMONLINE'
SCREENING COMPLETE - 0 REACTIONS TO VERIFY FROM 0 DOCUMENTS

100.0% DONE 0 VERIFIED 0 HIT RXNS 0 DOCS
SEARCH TIME: 00.00.10

3 FILES SEARCHED...

FULL SEARCH INITIATED 07:04:35 FILE 'PS'
SCREENING COMPLETE - 0 REACTIONS TO VERIFY FROM 0 DOCUMENTS

100.0% DONE 0 VERIFIED 0 HIT RXNS 0 DOCS
SEARCH TIME: 00.00.03

L5 0 L4

=> s l4 ful

FULL SEARCH INITIATED 07:04:48 FILE 'CASREACT'
SCREENING COMPLETE - 153 REACTIONS TO VERIFY FROM 36 DOCUMENTS

100.0% DONE 153 VERIFIED 0 HIT RXNS 0 DOCS
SEARCH TIME: 00.00.01

FULL SEARCH INITIATED 07:04:49 FILE 'CHEMINFORMRX'
SCREENING COMPLETE - 16 REACTIONS TO VERIFY FROM 8 DOCUMENTS

100.0% DONE 16 VERIFIED 0 HIT RXNS 0 DOCS
SEARCH TIME: 00.00.12

FULL SEARCH INITIATED 07:05:02 FILE 'DJSMONLINE'
SCREENING COMPLETE - 0 REACTIONS TO VERIFY FROM 0 DOCUMENTS

100.0% DONE 0 VERIFIED 0 HIT RXNS 0 DOCS
SEARCH TIME: 00.00.02

3 FILES SEARCHED...

FULL SEARCH INITIATED 07:05:05 FILE 'PS'
SCREENING COMPLETE - 0 REACTIONS TO VERIFY FROM 0 DOCUMENTS

100.0% DONE 0 VERIFIED 0 HIT RXNS 0 DOCS
SEARCH TIME: 00.00.01

L6 0 L4

=> file stnguide

COST IN U.S. DOLLARS SINCE FILE TOTAL

10/630,697

	ENTRY	SESSION
FULL ESTIMATED COST	362.09	363.18

FILE 'STNGUIDE' ENTERED AT 07:05:18 ON 31 AUG 2006
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	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	0.18	363.36

FILE 'REGISTRY' ENTERED AT 07:07:06 ON 31 AUG 2006
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=>Testing the current file.... screen

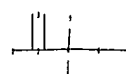
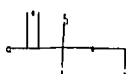
ENTER SCREEN EXPRESSION OR (END):end

=> screen 1992 OR 2016 OR 2021 OR 2026 OR 1838

L7 SCREEN CREATED

=>
Uploading C:\Documents and Settings\rkeys\My
Documents\STNEXP4\TEMPLATE\STANDARD\10630697a.str

10/630,697



chain nodes :
1 2 3 4 5 6 7 9 10
chain bonds :
1-2 2-3 2-4 4-5 4-6 4-9 6-7 7-10
exact/norm bonds :
2-3 4-6 4-9 6-7
exact bonds :
1-2 2-4 4-5 7-10

G1:F,CF3

Match level :
1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 9:CLASS 10:CLASS

L8 STRUCTURE UPLOADED

=> que L8 NOT L7

L9 QUE L8 NOT L7

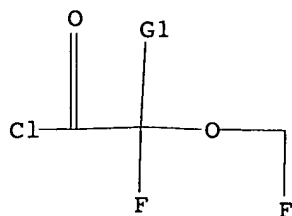
=> d

L9 HAS NO ANSWERS

L7 SCR 1992 OR 2016 OR 2021 OR 2026 OR 1838

L8 STR

10/630,697



G1 F,CF3

Structure attributes must be viewed using STN Express query preparation.
L9 QUE L8 NOT L7

=> s 19

SAMPLE SEARCH INITIATED 07:07:24 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 224 TO ITERATE

100.0% PROCESSED 224 ITERATIONS 4 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 3583 TO 5377
PROJECTED ANSWERS: 4 TO 200

L10 4 SEA SSS SAM L8 NOT L7

=> s 19 ful

FULL SEARCH INITIATED 07:07:33 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 4079 TO ITERATE

100.0% PROCESSED 4079 ITERATIONS 35 ANSWERS
SEARCH TIME: 00.00.01

L11 35 SEA SSS FUL L8 NOT L7

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	166.94	530.30

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FILE COVERS 1907 - 31 Aug 2006 VOL 145 ISS 10
FILE LAST UPDATED: 29 Aug 2006 (20060829/ED)

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soluble in both fluorous solvents and liquid carbon dioxide due to the attachment of a perfluoropolyether. The enzyme horse liver alc. dehydrogenase (HLADH) was active in catalyzing oxidation/reduction reactions using FNAD as a soluble coenzyme in a fluorous solvent, methoxynonafluorobutane (HFE), and liquid carbon dioxide. In both solvents, the activity of HLADH using FNAD was greater than the same molar amount of unmodified (insol.) NAD, indicating that a soluble coenzyme results in more efficient reactions.

RE.CNT 43 THERE ARE 43 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L15 ANSWER 2 OF 8 CAPLUS COPYRIGHT 2006 ACS on STN

AN 1993:625594 CAPLUS

DN 119:225594

TI Preparation of perfluorooxaalkanoyl halides and bis(perfluorooxaalkanoyl) peroxides

IN Sawada, Hideo; Matsumoto, Takeo; Nakayama, Masaharu

PA Nippon Oils & Fats Co Ltd, Japan

SO Jpn. Kokai Tokkyo Koho, 8 pp.

CODEN: JKXXAF

DT Patent

LA Japanese

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	JP 05170731	A2	19930709	JP 1991-341515	19911224
PRAI	JP 1991-341515		19911224		
AB	RfCOX and RfCO ₂ OCORf1 [Rf, Rf1 = CF ₃ [OCF(CF ₃)CF ₂] _n (OCF ₂) _m]; X = Br, Cl, F; n, m = 1-10] are prepared Chlorination of CF ₃ OCF(CF ₃)CF ₂ OCF ₂ CO ₂ H with POCl ₃ in DMF at 100° for 5 h gave 88% CF ₃ OCF(CF ₃)CF ₂ OCF ₂ COCl, which was treated with H ₂ O ₂ in CF ₃ CF ₂ CHCl ₂ -CCl ₂ CF ₂ CHFC1 mixture at temperature between -5° and +5° to give 79% [CF ₃ OCF(CF ₃)CF ₂ OCF ₂ CO ₂] ₂ .				

L15 ANSWER 3 OF 8 CAPLUS COPYRIGHT 2006 ACS on STN

AN 1986:625978 CAPLUS

DN 105:225978

TI Asymmetric addition of hydrogen cyanide to substituted benzaldehydes catalyzed by a synthetic cyclic peptide, cyclo[(S)-phenylalanyl-(S)-histidyl]

AU Kobayashi, Yoshiyuki; Asada, Shoichi; Watanabe, Ichigen; Hayashi, Hiroaki; Motoo, Yoshiyuki; Inoue, Shohei

CS Fac. Eng., Univ. Tokyo, Tokyo, 113, Japan

SO Bulletin of the Chemical Society of Japan (1986), 59(3), 893-5

CODEN: BCSJA8; ISSN: 0009-2673

DT Journal

LA English

OS CASREACT 105:225978

AB Optically active RC₆H₄CH(OH)CN (R = H, 4-Me, 3-Me, 2-Me, 3-MeO, 3-PhO) were prepared in 33-90% enantiomeric excess by addition of HCN to RC₆H₄CHO in C₆H₆ in the presence of cyclo[(S)-phenylalanyl-(S)-histidyl]. Highest optical yields were realized in nonpolar solvents, whereas, no asym. induction occurred in MeOH or Me₂SO.

L15 ANSWER 4 OF 8 CAPLUS COPYRIGHT 2006 ACS on STN

AN 1984:103740 CAPLUS

DN 100:103740

TI Synthesis of chiral steroid CD-ring synthon from D-leucine by means of diastereotopic face selection

AU Takahashi, Takashi; Okumoto, Hiroshi; Tsuji, Jiro; Harada, Nobuyuki

CS Dep. Chem. Eng., Tokyo Inst. Technol., Tokyo, 152, Japan

SO Journal of Organic Chemistry (1984), 49(5), 948-50

CODEN: JOCEAH; ISSN: 0022-3263

- DT Journal
LA English
AB The synthesis of indanone I is described, in which the correct absolute configuration is produced from D-leucine via the cis-vinyl iodide II (R = H), which serves to control the chirality of the rest of I by means of a remarkably effective diastereotopic face-selection. The optical purity of I and II (R = H) were checked by the ^{19}F NMR of their $(\text{F}_3\text{C})\text{CFOCF}(\text{CF}_3)\text{CO}_2\text{H}$ esters. The relative and absolute configuration of I and (-)-dienone III were determined by NMR and CD data. Thus, successive treatment of II (R = MeOCMe_2) with BuLi, CuI-PBu₃, 2-methyl-2-cyclopentenone, $\text{H}_2\text{C}:\text{C}(\text{SiMe}_3)\text{COMe}$, NaOMe, and HCl gave 58% I.
- L15 ANSWER 5 OF 8 CAPLUS COPYRIGHT 2006 ACS on STN
AN 1985:5467 CAPLUS
DN 102:5467
TI Thermal decomposition of (trifluoromethoxy)difluoroacetyl peroxide in heptane as a method of generating (trifluoromethoxy)difluoromethyl radicals
AU Komendantov, A. M.; Berenblit, V. V.; Sass, V.; Sokolov, S. V.
CS Vses. Nauchno-Issled. Inst. Sint. Kauch., USSR
SO Zhurnal Vsesoyuznogo Khimicheskogo Obshchestva im. D. I. Mendeleeva (1984), 29(3), 353-4
CODEN: ZVKOA6; ISSN: 0373-0247
DT Journal
LA Russian
AB The thermolysis of $(\text{CF}_3\text{OCF}_2\text{CO})_2\text{O}_2$ at 5-30° was a monomol., 1st-order process with activation energy 97 kJ/mol. The resulting $\text{CF}_3\text{OCF}_2\cdot$ radical abstracted H from heptane to give $\text{CF}_3\text{OCF}_2\text{H}$ quant. within 20 min at 50° with no side reaction.
- L15 ANSWER 6 OF 8 CAPLUS COPYRIGHT 2006 ACS on STN
AN 1984:422810 CAPLUS
DN 101:22810
TI Thermal decomposition of 2-(trifluoromethoxy)perfluoropropionyl peroxide in heptane as a method for the generation of 1-(trifluoromethoxy)perfluoroethyl radicals
AU Komendantov, A. M.; Starobin, Yu. K.; Berenblit, V. V.; Sass, V. P.; Sokolov, S. V.
CS Vses. Nauchno-Issled. Inst. Sint. Kauch., Leningrad, USSR
SO Zhurnal Vsesoyuznogo Khimicheskogo Obshchestva im. D. I. Mendeleeva (1984), 29(1), 113-14
CODEN: ZVKOA6; ISSN: 0373-0247
DT Journal
LA Russian
OS CASREACT 101:22810
AB The title thermolysis at 5-25° was 1st order in peroxide. The resulting $\cdot\text{CF}(\text{CF}_3)\text{OCF}_3$ abstracted H from the solvent to give $\text{CF}_3\text{OCHFOCF}_3$.
- L15 ANSWER 7 OF 8 CAPLUS COPYRIGHT 2006 ACS on STN
AN 1976:432486 CAPLUS
DN 85:32486
TI The reduction of perfluoroacyl halides with organosilicon hydrides. A direct synthesis of fluorine-containing esters and lactones
AU Croft, Thomas S.; McBrady, John J.
CS Cent. Res. Lab., 3M Co., St. Paul, MN, USA
SO Journal of Organic Chemistry (1976), 41(13), 2256-8
CODEN: JOCEAH; ISSN: 0022-3263
DT Journal
LA English
AB Reaction of perfluoroacyl halides with organosilicon halides in the presence of KF, ZnCl₂, and Pt/C gave 1,1-dihydroperfluoroalkyl perfluoroacylates. E.g., reaction of 2.8 g $\text{CF}_3\text{CF}_2\text{O}(\text{CF}_2)_2\text{COF}$ with Me_3SiH

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for 18 hr at 180° in the presence of the above catalysts gave 0.5 g CF₃CF₂O(CF₂)₂CO₂CH₂(CF₂)₂OCF₂CF₃. Similarly, I was prepared from perfluorocyclohexanecarbonyl fluoride; II (x = bond, CF₂, O) were prepared from perfluorosuccinyl or perfluoroglutaryl fluoride or O(CF₂COCl)₂.

L15 ANSWER 8 OF 8 CAPLUS COPYRIGHT 2006 ACS on STN

AN 1975:594115 CAPLUS

DN 83:194115

TI Perfluorinated linear polyethers having reactive terminal groups at both ends of the chain

IN Sianesi, Dario; Caporiccio, Gerardo; Mensi, Domenico

PA Montedison S.p.A., Italy

SO U.S., 14 pp.

CODEN: USXXAM

DT Patent

LA English

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 3847978	A	19741112	US 1969-834486	19690618
PRAI	US 1968-787309	A2	19681226		

AB Perfluorinated linear polyethers containing peroxidic linkages were chain-cleaved by reducing agents to give bifunctional perfluorinated linear oligopolyethers with chemical-reactive terminal groups. Thus, hexafluoropropene [116-15-4] was treated with oxygen under the influence of uv light to give a peroxidized poly(perfluoropropylene oxide) [25038-02-2] which was reduced by H over a Pd catalyst to give a series of carboxy- and trifluoroacetyl-terminated oligopolyethers. One of these, CF₃COCF₂O(C₃F₆O)₂CF(CF₃)CO₂H [42775-40-6], boiling point 210-2°, formed a polymer with hexamethylenediamine [55809-69-3].

=> file stnguide

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
44.07	574.37

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
-6.00	-6.00

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LAST RELOADED: Aug 25, 2006 (20060825/UP).

=> log y

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
0.42	574.79

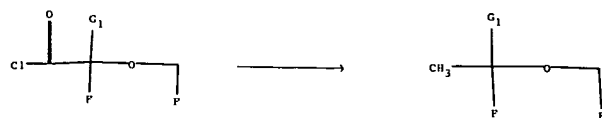
FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
0.00	-6.00

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STN INTERNATIONAL LOGOFF AT 07:17:27 ON 31 AUG 2006



chain nodes :

1 2 3 4 5 6 7 9 10 11 12 13 15 16 17 18

chain bonds :

1-2 2-3 2-4 4-5 4-6 4-9 6-7 7-18 10-11 11-12 11-13 11-15 15-16 16-17

exact/norm bonds :

2-3 4-6 4-9 6-7 11-13 11-15 15-16

exact bonds :

1-2 2-4 4-5 7-18 10-11 11-12 16-17

G1:F,CF3

Match level :

1:CLASS2:CLASS3:CLASS4:CLASS5:CLASS6:CLASS7:CLASS9:CLASS10:CLASS11:CLASS12:CLASS13:CLASS15:CLASS16:CLASS17:CLASS18:CLASS

fragments assigned product role:

containing 10

fragments assigned reactant/reagent role:

containing 1